Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070316\
Data File : BF088674.D
Acq On : 4 Jul 2016 4:11
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 04 07:01:42 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 30 18:01:55 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	93	0.01
2		1,4-Dioxane	40.000	35.911	10.2	86	-0.01
3		Pyridine	40.000	36.320	9.2	88	0.00
4		n-Nitrosodimethylamine	40.000	34.375	14.1	83	-0.01
5	S	2-Fluorophenol	80.000	70.980	11.3	82	0.00
6		Aniline	40.000	35.679	10.8	83	0.01
7	S	Phenol-d6	80.000	72.819	9.0	89	0.01
8		2-Chlorophenol	40.000	40.148	-0.4	101	0.01
9		Benzaldehyde	40.000	35.085	12.3	88	0.01
10	C	Phenol	40.000	36.613	8.5	85	0.00
11		bis(2-Chloroethyl)ether	40.000	37.524	6.2	93	0.01
12		1,3-Dichlorobenzene	40.000	37.684	5.8	96	0.01
13	C	1,4-Dichlorobenzene	40.000	36.564	8.6	91	0.00
14		1,2-Dichlorobenzene	40.000	40.444	-1.1	105	0.01
15		Benzyl Alcohol	40.000	36.562	8.6	83	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	34.908	12.7	83	0.01
17		2-Methylphenol	40.000	39.957	0.1	93	0.01
18		Hexachloroethane	40.000	39.550	1.1	94	0.01
19	P	n-Nitroso-di-n-propylamine	40.000	38.670	3.3	105	0.01
20		3+4-Methylphenols	40.000	34.568	13.6	86	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	89	0.00
22		Acetophenone	40.000	37.302	6.7	85	0.00
23	S	Nitrobenzene-d5	80.000	81.560	-2.0	96	0.00
24		Nitrobenzene	40.000	41.272	-3.2	96	0.01
25		Isophorone	40.000	38.146	4.6	89	0.00
26	C	2-Nitrophenol	40.000	48.150	-20.4#	116	0.01
27		2,4-Dimethylphenol	40.000	41.250	-3.1	90	0.01
28		bis(2-Chloroethoxy)methane	40.000	38.870	2.8	94	0.01
29	C	2,4-Dichlorophenol	40.000	42.357	-5.9	100	0.01
30		1,2,4-Trichlorobenzene	40.000	39.735	0.7	89	0.00
31		Naphthalene	40.000	37.994	5.0	84	0.00
32		Benzoic acid	40.000	39.920	0.2	90	0.00
33		4-Chloroaniline	40.000	44.261	-10.7	100	0.01
34	C	Hexachlorobutadiene	40.000	42.076	-5.2	94	0.01
35		Caprolactam	40.000	40.621	-1.6	88	0.00
36	C	4-Chloro-3-methylphenol	40.000	44.207	-10.5	100	0.01
37		2-Methylnaphthalene	40.000	41.648	-4.1	105	0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	92	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.540	3.7	86	0.00
40	P	Hexachlorocyclopentadiene	40.000	36.297	9.3	85	0.00
41	S	2,4,6-Tribromophenol	80.000	80.730	-0.9	89	0.00
42	C	2,4,6-Trichlorophenol	40.000	43.123	-7.8	109	0.01
43		2,4,5-Trichlorophenol	40.000	41.794	-4.5	94	0.00
44	S	2-Fluorobiphenyl	80.000	88.744	-10.9	113	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070316\
Data File : BF088674.D
Acq On : 4 Jul 2016 4:11
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 04 07:01:42 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 30 18:01:55 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	38.014	5.0	86	0.00
46		2-Chloronaphthalene	40.000	38.731	3.2	87	0.00
47		2-Nitroaniline	40.000	37.464	6.3	78	0.00
48		Acenaphthylene	40.000	38.002	5.0	87	0.00
49		Dimethylphthalate	40.000	44.399	-11.0	111	0.01
50		2,6-Dinitrotoluene	40.000	47.959	-19.9	119	0.01
51	C	Acenaphthene	40.000	40.605	-1.5	98	0.00
52		3-Nitroaniline	40.000	37.644	5.9	77	0.00
53	P	2,4-Dinitrophenol	40.000	39.314	1.7	93	0.00
54		Dibenzofuran	40.000	38.178	4.6	89	0.00
55	P	4-Nitrophenol	40.000	37.312	6.7	78	0.00
56		2,4-Dinitrotoluene	40.000	49.058	-22.6	119	0.01
57		Fluorene	40.000	36.358	9.1	81	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	44.283	-10.7	99	0.01
59		Diethylphthalate	40.000	43.055	-7.6	100	0.01
60		4-Chlorophenyl-phenylether	40.000	42.983	-7.5	108	0.01
61		4-Nitroaniline	40.000	45.920	-14.8	115	0.01
62		Azobenzene	40.000	41.483	-3.7	100	0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	99	0.01
64		4,6-Dinitro-2-methylphenol	40.000	36.587	8.5	82	0.00
65	c	n-Nitrosodiphenylamine	40.000	36.343	9.1	87	0.00
66		4-Bromophenyl-phenylether	40.000	37.854	5.4	96	0.00
67		Hexachlorobenzene	40.000	40.401	-1.0	100	0.01
68		Atrazine	40.000	35.142	12.1	84	0.00
69	C	Pentachlorophenol	40.000	41.779	-4.4	98	0.01
70		Phenanthrene	40.000	40.457	-1.1	106	0.01
71		Anthracene	40.000	35.334	11.7	88	0.00
72		Carbazole	40.000	40.705	-1.8	112	0.01
73		Di-n-butylphthalate	40.000	38.667	3.3	101	0.00
74	C	Fluoranthene	40.000	36.759	8.1	87	0.01
75	I	Chrysene-d12	20.000	20.000	0.0	87	0.00
76		Benzidine	40.000	40.433	-1.1	85	0.01
77		Pyrene	40.000	42.689	-6.7	88	0.01
78	S	Terphenyl-d14	80.000	80.994	-1.2	90	0.01
79		Butylbenzylphthalate	40.000	44.625	-11.6	97	0.01
80		Benzo(a)anthracene	40.000	39.653	0.9	86	0.00
81		3,3'-Dichlorobenzidine	40.000	40.574	-1.4	91	0.01
82		Chrysene	40.000	42.242	-5.6	93	0.01
83		Bis(2-ethylhexyl)phthalate	40.000	46.083	-15.2	95	0.01
84	c	Di-n-octyl phthalate	40.000	45.964	-14.9	95	0.01
85		Indeno(1,2,3-cd)pyrene	40.000	40.159	-0.4	91	0.00
86	I	Perylene-d12	20.000	20.000	0.0	83	0.01
87		Benzo(b)fluoranthene	40.000	40.811	-2.0	89	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070316\
Data File : BF088674.D
Acq On : 4 Jul 2016 4:11
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 04 07:01:42 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 30 18:01:55 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	35.058	12.4	73	0.00
89 C	Benzo(a)pyrene	40.000	39.227	1.9	81	0.01
90	Dibenzo(a,h)anthracene	40.000	43.007	-7.5	91	0.01
91	Benzo(g,h,i)perylene	40.000	46.154	-15.4	98	0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 1